



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:38 PM UTC

PDB ID : 3DK5 / pdb_00003dk5
Title : Crystal Structure of Apo-GlmU from Mycobacterium tuberculosis
Authors : Verma, S.K.; Prakash, B.
Deposited on : 2008-06-24
Resolution : 2.23 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

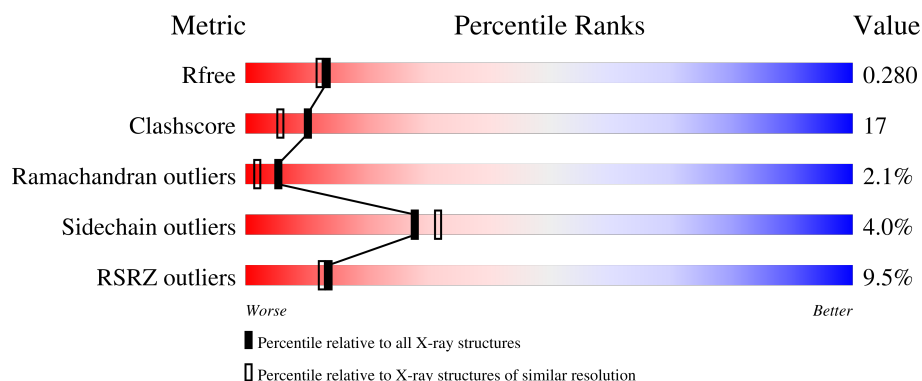
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.23 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3416 (2.26-2.22)
Clashscore	190562	3556 (2.26-2.22)
Ramachandran outliers	187476	3500 (2.26-2.22)
Sidechain outliers	187428	3501 (2.26-2.22)
RSRZ outliers	180081	3415 (2.26-2.22)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	495	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3248 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Bifunctional protein glmU.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	442	Total	C	N	O	S	0	0	0
			3144	1950	567	622	5			

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mg	0	0
			2	2		

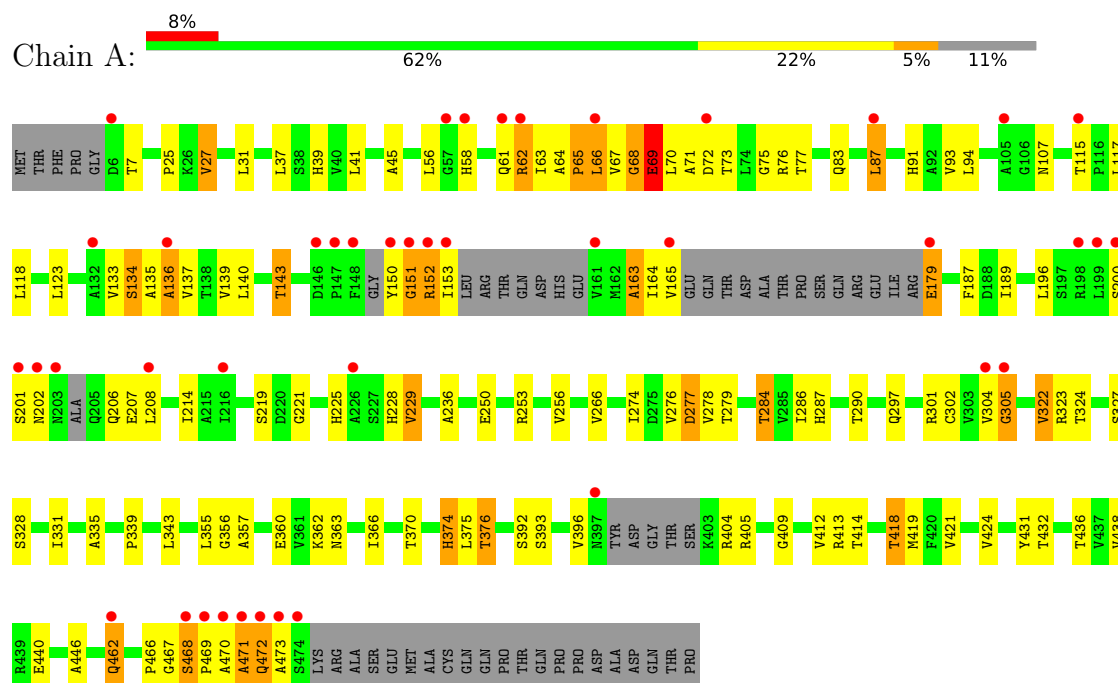
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	102	Total	O	0	0
			102	102		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Bifunctional protein glmU



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	79.60Å 79.60Å 278.00Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.30 – 2.23 29.30 – 2.23	Depositor EDS
% Data completeness (in resolution range)	100.0 (29.30-2.23) 98.1 (29.30-2.23)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.63 (at 2.24Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.225 , 0.271 0.241 , 0.280	Depositor DCC
R_{free} test set	1567 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	32.2	Xtriage
Anisotropy	0.343	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 41.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.048 for -h-k,k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3248	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.72% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.61	29/3188 (0.9%)	1.35	12/4363 (0.3%)

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	343	LEU	CA-C	8.37	1.63	1.52
1	A	229	VAL	CA-CB	7.01	1.62	1.53
1	A	360	GLU	CA-C	6.91	1.61	1.52
1	A	322	VAL	CA-C	6.82	1.60	1.52
1	A	335	ALA	CA-CB	-6.80	1.42	1.53
1	A	256	VAL	CA-CB	6.69	1.61	1.54
1	A	297	GLN	N-CA	6.18	1.53	1.46
1	A	409	GLY	N-CA	5.97	1.52	1.45
1	A	324	THR	CA-C	5.96	1.59	1.52
1	A	266	VAL	N-CA	5.88	1.53	1.46
1	A	366	ILE	CA-CB	5.79	1.61	1.54
1	A	355	LEU	N-CA	5.78	1.53	1.46
1	A	446	ALA	CA-CB	-5.69	1.44	1.53
1	A	370	THR	CA-C	5.62	1.60	1.52
1	A	274	ILE	CA-C	5.62	1.60	1.52
1	A	279	THR	N-CA	5.55	1.52	1.46
1	A	45	ALA	CA-CB	-5.55	1.44	1.53
1	A	418	THR	CA-CB	5.52	1.61	1.53
1	A	290	THR	C-O	-5.32	1.17	1.24
1	A	117	LEU	C-O	5.32	1.30	1.23
1	A	93	VAL	CA-CB	5.31	1.61	1.54
1	A	466	PRO	CA-C	5.31	1.59	1.52
1	A	421	VAL	CA-CB	-5.25	1.48	1.53
1	A	412	VAL	CA-CB	-5.18	1.47	1.54
1	A	374	HIS	N-CA	5.06	1.52	1.46
1	A	286	ILE	CA-CB	5.06	1.60	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	419	MET	N-CA	-5.05	1.40	1.46
1	A	331	ILE	C-O	-5.03	1.18	1.24
1	A	356	GLY	N-CA	5.02	1.51	1.45

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	202	ASN	N-CA-C	7.45	120.65	109.25
1	A	305	GLY	CA-C-N	-6.79	111.35	119.84
1	A	305	GLY	C-N-CA	-6.79	111.35	119.84
1	A	87	LEU	N-CA-C	6.29	119.12	111.82
1	A	201	SER	N-CA-C	5.92	123.41	110.80
1	A	396	VAL	CB-CA-C	-5.64	103.27	111.68
1	A	301	ARG	N-CA-C	5.54	119.33	112.47
1	A	65	PRO	N-CA-C	5.49	120.98	113.84
1	A	206	GLN	N-CA-C	5.34	119.19	111.56
1	A	68	GLY	N-CA-C	-5.26	100.71	113.18
1	A	163	ALA	N-CA-C	5.21	115.42	108.23
1	A	56	LEU	N-CA-C	5.02	117.29	109.52

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3144	0	3101	107	0
2	A	2	0	0	0	0
3	A	102	0	0	0	0
All	All	3248	0	3101	107	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 17.

All (107) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:ARG:O	1:A:65:PRO:HD2	1.37	1.25
1:A:468:SER:OG	1:A:469:PRO:HD3	1.36	1.22
1:A:143:THR:HG22	1:A:228:HIS:HE1	1.05	1.10
1:A:143:THR:HG22	1:A:228:HIS:CE1	1.86	1.09
1:A:468:SER:CB	1:A:469:PRO:CD	2.32	1.08
1:A:468:SER:HB3	1:A:469:PRO:HD2	1.37	1.03
1:A:276:VAL:O	1:A:277:ASP:HB2	1.62	0.99
1:A:200:SER:O	1:A:208:LEU:HG	1.60	0.99
1:A:468:SER:HB3	1:A:469:PRO:CD	1.94	0.96
1:A:468:SER:OG	1:A:469:PRO:CD	2.14	0.94
1:A:83:GLN:HE21	1:A:91:HIS:HD2	1.20	0.89
1:A:25:PRO:HB3	1:A:58:HIS:CE1	2.10	0.87
1:A:467:GLY:O	1:A:468:SER:HB3	1.72	0.86
1:A:62:ARG:O	1:A:65:PRO:CD	2.23	0.86
1:A:68:GLY:O	1:A:69:GLU:HB2	1.76	0.86
1:A:25:PRO:HG3	1:A:58:HIS:NE2	1.92	0.84
1:A:472:GLN:O	1:A:473:ALA:HB3	1.77	0.84
1:A:135:ALA:O	1:A:136:ALA:HB3	1.80	0.82
1:A:468:SER:CB	1:A:469:PRO:HD3	2.04	0.82
1:A:468:SER:HG	1:A:469:PRO:HD3	1.42	0.82
1:A:374:HIS:O	1:A:376:THR:HG23	1.80	0.81
1:A:150:TYR:O	1:A:151:GLY:O	1.99	0.80
1:A:467:GLY:O	1:A:468:SER:CB	2.28	0.79
1:A:37:LEU:HD21	1:A:66:LEU:HD21	1.65	0.79
1:A:200:SER:O	1:A:208:LEU:CG	2.30	0.78
1:A:83:GLN:HE21	1:A:91:HIS:CD2	2.04	0.75
1:A:150:TYR:O	1:A:151:GLY:C	2.31	0.74
1:A:67:VAL:CA	1:A:71:ALA:HB3	2.19	0.73
1:A:27:VAL:HG13	1:A:37:LEU:HB2	1.69	0.73
1:A:64:ALA:HB3	1:A:65:PRO:HD3	1.72	0.70
1:A:67:VAL:CA	1:A:71:ALA:H	2.05	0.70
1:A:135:ALA:O	1:A:136:ALA:CB	2.39	0.69
1:A:276:VAL:O	1:A:277:ASP:CB	2.38	0.69
1:A:67:VAL:CA	1:A:71:ALA:CB	2.70	0.68
1:A:143:THR:CG2	1:A:228:HIS:CE1	2.72	0.68
1:A:472:GLN:O	1:A:473:ALA:CB	2.39	0.67
1:A:143:THR:CG2	1:A:228:HIS:HE1	1.95	0.67
1:A:25:PRO:CG	1:A:58:HIS:NE2	2.58	0.66
1:A:470:ALA:O	1:A:471:ALA:HB3	1.95	0.66
1:A:7:THR:HG22	1:A:107:ASN:HB2	1.77	0.66
1:A:58:HIS:HE1	1:A:62:ARG:NE	1.94	0.64
1:A:62:ARG:C	1:A:65:PRO:HD2	2.20	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:HIS:CG	1:A:58:HIS:O	2.51	0.64
1:A:25:PRO:HB3	1:A:58:HIS:HE1	1.60	0.62
1:A:71:ALA:HB1	1:A:76:ARG:O	1.99	0.62
1:A:58:HIS:HE1	1:A:62:ARG:CZ	2.13	0.61
1:A:133:VAL:HG21	1:A:225:HIS:CD2	2.36	0.60
1:A:58:HIS:CE1	1:A:62:ARG:NE	2.70	0.60
1:A:374:HIS:O	1:A:376:THR:CG2	2.49	0.59
1:A:471:ALA:N	1:A:472:GLN:O	2.33	0.59
1:A:200:SER:O	1:A:208:LEU:CD1	2.51	0.58
1:A:468:SER:O	1:A:469:PRO:C	2.46	0.58
1:A:66:LEU:O	1:A:68:GLY:N	2.36	0.58
1:A:133:VAL:O	1:A:134:SER:O	2.22	0.57
1:A:470:ALA:O	1:A:471:ALA:CB	2.52	0.56
1:A:63:ILE:HA	1:A:66:LEU:HD12	1.87	0.56
1:A:133:VAL:HG21	1:A:225:HIS:HD2	1.71	0.56
1:A:41:LEU:HD12	1:A:70:LEU:HD13	1.89	0.55
1:A:287:HIS:HB2	1:A:305:GLY:C	2.32	0.55
1:A:94:LEU:HD13	1:A:196:LEU:HD13	1.90	0.54
1:A:72:ASP:O	1:A:75:GLY:N	2.41	0.53
1:A:327:SER:O	1:A:328:SER:C	2.54	0.51
1:A:322:VAL:O	1:A:323:ARG:C	2.53	0.51
1:A:153:ILE:HG23	1:A:153:ILE:O	2.11	0.50
1:A:76:ARG:NH1	1:A:77:THR:O	2.45	0.50
1:A:136:ALA:O	1:A:187:PHE:HA	2.12	0.50
1:A:418:THR:HA	1:A:436:THR:O	2.12	0.50
1:A:375:LEU:O	1:A:392:SER:HA	2.13	0.49
1:A:76:ARG:HH11	1:A:76:ARG:HG3	1.77	0.49
1:A:87:LEU:O	1:A:207:GLU:HB3	2.12	0.49
1:A:404:ARG:O	1:A:405:ARG:HB2	2.12	0.49
1:A:287:HIS:HB2	1:A:305:GLY:O	2.14	0.48
1:A:339:PRO:O	1:A:357:ALA:HA	2.13	0.48
1:A:115:THR:HG23	1:A:236:ALA:O	2.14	0.47
1:A:376:THR:HG22	1:A:393:SER:H	1.79	0.47
1:A:72:ASP:O	1:A:73:THR:C	2.58	0.47
1:A:153:ILE:HG22	1:A:179:GLU:O	2.15	0.46
1:A:25:PRO:CB	1:A:62:ARG:HH21	2.28	0.46
1:A:153:ILE:HD13	1:A:179:GLU:OE1	2.15	0.46
1:A:25:PRO:CB	1:A:58:HIS:CE1	2.91	0.46
1:A:462:GLN:NE2	1:A:473:ALA:HB1	2.31	0.46
1:A:277:ASP:O	1:A:278:VAL:C	2.59	0.45
1:A:68:GLY:O	1:A:69:GLU:CB	2.54	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:362:LYS:O	1:A:363:ASN:C	2.60	0.45
1:A:414:THR:HA	1:A:432:THR:HB	1.98	0.45
1:A:304:VAL:O	1:A:305:GLY:O	2.35	0.44
1:A:163:ALA:C	1:A:164:ILE:HG13	2.42	0.44
1:A:58:HIS:CE1	1:A:62:ARG:CZ	2.99	0.43
1:A:250:GLU:HA	1:A:253:ARG:NH1	2.33	0.43
1:A:140:LEU:HD22	1:A:229:VAL:HG23	2.00	0.43
1:A:413:ARG:O	1:A:431:TYR:HA	2.18	0.43
1:A:424:VAL:HG21	1:A:438:VAL:HG12	2.01	0.43
1:A:133:VAL:CG2	1:A:225:HIS:HD2	2.31	0.42
1:A:137:VAL:HG21	1:A:214:ILE:CD1	2.49	0.42
1:A:284:THR:HG22	1:A:302:CYS:H	1.83	0.42
1:A:133:VAL:O	1:A:134:SER:C	2.61	0.42
1:A:219:SER:C	1:A:221:GLY:H	2.26	0.42
1:A:58:HIS:O	1:A:58:HIS:ND1	2.53	0.42
1:A:468:SER:O	1:A:471:ALA:O	2.38	0.41
1:A:151:GLY:O	1:A:152:ARG:CB	2.68	0.41
1:A:376:THR:HG22	1:A:392:SER:N	2.36	0.41
1:A:118:LEU:HD21	1:A:123:LEU:HD21	2.02	0.41
1:A:61:GLN:O	1:A:65:PRO:HD3	2.20	0.41
1:A:322:VAL:HG12	1:A:323:ARG:HG3	2.03	0.40
1:A:31:LEU:O	1:A:39:HIS:HE1	2.05	0.40
1:A:467:GLY:H	1:A:470:ALA:HB3	1.86	0.40
1:A:62:ARG:O	1:A:63:ILE:C	2.62	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	430/495 (87%)	387 (90%)	34 (8%)	9 (2%)	5 2

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	151	GLY
1	A	468	SER
1	A	134	SER
1	A	136	ALA
1	A	277	ASP
1	A	69	GLU
1	A	152	ARG
1	A	471	ALA
1	A	472	GLN

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	327/388 (84%)	314 (96%)	13 (4%)	28	32

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	VAL
1	A	62	ARG
1	A	66	LEU
1	A	69	GLU
1	A	139	VAL
1	A	143	THR
1	A	165	VAL
1	A	179	GLU
1	A	189	ILE
1	A	284	THR
1	A	376	THR
1	A	440	GLU
1	A	462	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	HIS
1	A	58	HIS
1	A	91	HIS
1	A	228	HIS
1	A	363	ASN
1	A	454	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	442/495 (89%)	0.35	42 (9%) 14 12	17, 39, 74, 97	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	151	GLY	7.0
1	A	165	VAL	6.6
1	A	305	GLY	6.4
1	A	150	TYR	6.0
1	A	58	HIS	6.0
1	A	471	ALA	5.8
1	A	62	ARG	5.3
1	A	472	GLN	5.3
1	A	470	ALA	4.8
1	A	469	PRO	4.5
1	A	153	ILE	4.2
1	A	226	ALA	3.7
1	A	200	SER	3.6
1	A	72	ASP	3.5
1	A	468	SER	3.4
1	A	474	SER	3.4
1	A	152	ARG	3.4
1	A	201	SER	3.4
1	A	202	ASN	3.3
1	A	87	LEU	3.2
1	A	208	LEU	2.9
1	A	304	VAL	2.9
1	A	148	PHE	2.9
1	A	473	ALA	2.8
1	A	179	GLU	2.6
1	A	198	ARG	2.6
1	A	146	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	6	ASP	2.5
1	A	66	LEU	2.5
1	A	136	ALA	2.5
1	A	57	GLY	2.5
1	A	199	LEU	2.4
1	A	161	VAL	2.4
1	A	397	ASN	2.2
1	A	147	PRO	2.2
1	A	61	GLN	2.2
1	A	115	THR	2.1
1	A	203	ASN	2.1
1	A	216	ILE	2.1
1	A	462	GLN	2.1
1	A	132	ALA	2.0
1	A	105	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MG	A	497	1/1	0.80	0.44	26,26,26,26	1
2	MG	A	496	1/1	0.93	0.28	17,17,17,17	1

6.5 Other polymers [i](#)

There are no such residues in this entry.